

V	Final Report
	Re-Issued Repor
	Revised Report
	port Date:

# Laboratory Report

Gulf Oil L.P. 281 Eastern Avenue Chelsea, MA 02150 Attn: Andrew P. Adams

Project: Gulf Terminal - Chelsea, MA

Project #: Gulf Chelsea

Laboratory IDClient Sample IDMatrixDate SampledDate ReceivedSC29001-01Outfall 003Surface Water30-Nov-16 07:0001-Dec-16 15:00

I attest that the information contained within the report has been reviewed for accuracy and checked against the quality control requirements for each method. These results relate only to the sample(s) as received.

All applicable NELAC requirements have been met.

Massachusetts # M-MA138/MA1110 Connecticut # PH-0777 Florida # E87936 Maine # MA138 New Hampshire # 2972/2538 New Jersey # MA011 New York # 11393 Pennsylvania # 68-04426/68-02924 Rhode Island # LAO00348 USDA # P330-15-00375 Vermont # VT-11393



Authorized by:

June O'Connor Laboratory Director

Eurofins Spectrum Analytical holds primary certification in the State of Massachusetts for the analytes as indicated with an X in the "Cert." column within this report. Please note that the State of Massachusetts does not offer certification for all analytes. Please refer to our website for specific certification holdings in each state.

Please note that this report contains 11 pages of analytical data plus Chain of Custody document(s). When the Laboratory Report is indicated as revised, this report supersedes any previously dated reports for the laboratory ID(s) referenced above. Where this report identifies subcontracted analyses, copies of the subcontractor's test report are available upon request. This report may not be reproduced, except in full, without written approval from Eurofins Spectrum Analytical, Inc.

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Please contact the Laboratory or Technical Director at 800-789-9115 with any questions regarding the data contained in this laboratory report.

### **CASE NARRATIVE:**

Data has been reported to the RDL. This report excludes estimated concentrations detected below the RDL and above the MDL (J-Flag).

All non-detects and all results below the reporting limit are reported as "<" (less than) the reporting limit in this report.

The samples were received 3.0 degrees Celsius, please refer to the Chain of Custody for details specific to temperature upon receipt. An infrared thermometer with a tolerance of  $\pm 1.0$  degrees Celsius was used immediately upon receipt of the samples.

If a Matrix Spike (MS), Matrix Spike Duplicate (MSD) or Duplicate (DUP) was not requested on the Chain of Custody, method criteria may have been fulfilled with a source sample not of this Sample Delivery Group.

Analyses for Total Hardness, pH, and Total Residual Chlorine fall under the state of Pennsylvania code Chapter 252.6 accreditation by rule.

See below for any non-conformances and issues relating to quality control samples and/or sample analysis/matrix.

## SW846 8260C

### Calibration:

1612010

Analyte quantified by quadratic equation type calibration.

Naphthalene

This affected the following samples:

1621081-BLK1 1621081-BS1 1621081-BSD1

Outfall 003

S610283-ICV1

S610389-CCV1

# **Sample Acceptance Check Form**

Client:

Gulf Oil L.P.

Project:	Gulf Terminal - Chelsea, MA / Gulf Chelsea			
Work Order:	SC29001			
Sample(s) received on:	12/1/2016			
The following outlines the	condition of samples for the attached Chain of Custody upon receipt.			
		Yes	<u>No</u>	<u>N/A</u>
Were custody seal	s present?		$\overline{\checkmark}$	
Were custody seal	s intact?			<b>✓</b>
Were samples rece	eived at a temperature of $\leq 6$ °C?	$\checkmark$		
Were samples refr	igerated upon transfer to laboratory representative?	$\checkmark$		
Were sample conta	ainers received intact?	$\checkmark$		
	perly labeled (labels affixed to sample containers and include sample ID, site roject number and the collection date)?	$\checkmark$		
Were samples acco	ompanied by a Chain of Custody document?	$\checkmark$		
include sample ID	stody document include proper, full, and complete documentation, which shall stody, site location, and/or project number, date and time of collection, collector's name, sample matrix and any special remarks concerning the sample?	<b>V</b>		
Did sample contai	ner labels agree with Chain of Custody document?	$\checkmark$		
Were samples rece	eived within method-specific holding times?	$\checkmark$		

# **Summary of Hits**

**Lab ID:** SC29001-01

Client ID: Outfall 003

Parameter	Result	Flag	Reporting Limit	Units	Analytical Method
Total Suspended Solids	10.3		0.5	mg/l	SM2540D (11)

Please note that because there are no reporting limits associated with hazardous waste characterizations or micro analyses, this summary does not include hits from these analyses if included in this work order.

Sample Id Outfall 00 SC29001-				lient Pro Gulf Cho			<u>Matrix</u> Surface Wa	· · · · · · · · · · · · · · · · · · ·	ection Date -Nov-16 07			ceived Dec-16	
CAS No.	Analyte(s)	Result	Flag U	nits	*RDL	MDL	Dilution	Method Ref.	Prepared	Analyzed	Analyst	Batch	Cert.
Volatile O	rganic Compounds												
-	rganic Aromatics by SW8 by method SW846 5030												
71-43-2	Benzene	< 1.0	ŀ	ug/l	1.0	0.3	1	SW846 8260C	02-Dec-16	03-Dec-16	TS	1621081	
91-20-3	Naphthalene	< 1.0	ŀ	ug/l	1.0	0.3	1	"	"	"	"	u u	
Surrogate i	recoveries:												
460-00-4	4-Bromofluorobenzene	78			70-13	0 %			"	"	"	"	
2037-26-5	Toluene-d8	102			70-13	0 %			"	"	"	"	
17060-07-0	1,2-Dichloroethane-d4	117			70-13	0 %			"	"	"	"	
1868-53-7	Dibromofluoromethane	115			70-13	0 %		"	"	"	"	"	
Semivolati	le Organic Compounds by	GCMS											
SVOCs by Prepared	<u>y SIM</u> by method SW846 35100	<u>2</u>											
50-32-8	Benzo (a) pyrene	< 0.051	ŀ	ıg/l	0.051	0.036	1	SW846 8270D SIM	05-Dec-16	06-Dec-16	MSL	1621185	
91-20-3	Naphthalene	< 0.051	ŀ	ug/l	0.051	0.027	1	"	п	"	"	"	
Surrogate i	recoveries:												
205440-82-0	Benzo (e) pyrene-d12	93			30-13	0 %		"	"	"	"	"	
General C	hemistry Parameters												
	рН	6.74	рН	Units			1	ASTM D 1293-99B	01-Dec-16 17:29	01-Dec-16 19:18	BD	1621045	Χ
	Total Suspended Solids	10.3	n	ng/l	0.5	0.2	1	SM2540D (11)	02-Dec-16	06-Dec-16	CMB	1621065	Χ
	cted Analyses by method 369621												
Analysis pe	erformed by Phoenix Environ	nmental Labs, Inc.	* - MACT007										
	Oil and Grease by EPA 1664A	< 1.4	m	ng/L	1.4	1.4	1	E1664A		12-Dec-16 08:26	MACT0	369621A	

# **Volatile Organic Compounds - Quality Control**

					Spike	Source		%REC		RPD
nalyte(s)	Result	Flag	Units	*RDL	Level	Result	%REC	Limits	RPD	Limi
atch 1621081 - SW846 5030 Water MS										
Blank (1621081-BLK1)					Pre	epared & A	nalyzed: 02-	-Dec-16		
Benzene	< 1.0		μg/l	1.0						
Ethylbenzene	< 1.0		μg/l	1.0						
Methyl tert-butyl ether	< 1.0		μg/l	1.0						
Naphthalene	< 1.0		μg/l	1.0						
Toluene	< 1.0		μg/l	1.0						
m,p-Xylene	< 2.0		μg/l	2.0						
o-Xylene	< 1.0		μg/l	1.0						
Surrogate: 4-Bromofluorobenzene	40.2		μg/l		50.0		80	70-130		
Surrogate: Toluene-d8	49.3		μg/l		50.0		99	70-130		
Surrogate: 1,2-Dichloroethane-d4	55.7		μg/l		50.0		111	70-130		
Surrogate: Dibromofluoromethane	53.0		μg/l		50.0		106	70-130		
LCS (1621081-BS1)					Pre	epared & A	nalyzed: 02-	-Dec-16		
Benzene	19.6		μg/l		20.0		98	70-130		
Ethylbenzene	17.9		μg/l		20.0		90	70-130		
Methyl tert-butyl ether	20.7		μg/l		20.0		104	70-130		
Naphthalene	19.3		μg/l		20.0		96	70-130		
Toluene	18.5		μg/l		20.0		92	70-130		
m,p-Xylene	17.8		μg/l		20.0		89	70-130		
o-Xylene	18.5		μg/l		20.0		92	70-130		
Surrogate: 4-Bromofluorobenzene	54.7		μg/l		50.0		109	70-130		
Surrogate: Toluene-d8	49.4		μg/l		50.0		99	70-130		
Surrogate: 1,2-Dichloroethane-d4	48.7		μg/l		50.0		97	70-130		
Surrogate: Dibromofluoromethane	48.4		μg/l		50.0		97	70-130		
LCS Dup (1621081-BSD1)					Pre	epared: 02-	Dec-16 An	alyzed: 03-D	ec-16	
Benzene	20.4		μg/l		20.0		102	70-130	4	20
Ethylbenzene	18.1		μg/l		20.0		90	70-130	1	20
Methyl tert-butyl ether	21.3		μg/l		20.0		106	70-130	3	20
Naphthalene	19.8		μg/l		20.0		99	70-130	3	20
Toluene	18.8		μg/l		20.0		94	70-130	2	20
m,p-Xylene	18.2		μg/l		20.0		91	70-130	2	20
o-Xylene	19.4		μg/l		20.0		97	70-130	5	20
Surrogate: 4-Bromofluorobenzene	54.3		μg/l		50.0		109	70-130		
Surrogate: Toluene-d8	49.4		μg/l		50.0		99	70-130		
Surrogate: 1,2-Dichloroethane-d4	48.8		μg/l		50.0		98	70-130		
Surrogate: Dibromofluoromethane	47.8		μg/l		50.0		96	70-130		

# Semivolatile Organic Compounds by GCMS - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Batch 1621185 - SW846 3510C										
Blank (1621185-BLK2)					Pre	epared: 05-	Dec-16 An	alyzed: 06-D	ec-16	
Acenaphthene	< 0.050		μg/l	0.050						
Acenaphthylene	< 0.050		μg/l	0.050						
Anthracene	< 0.050		μg/l	0.050						
Benzo (a) anthracene	< 0.050		μg/l	0.050						
Benzo (a) pyrene	< 0.050		μg/l	0.050						
Benzo (b) fluoranthene	< 0.050		μg/l	0.050						
Benzo (g,h,i) perylene	< 0.050		μg/l	0.050						
Benzo (k) fluoranthene	< 0.050		μg/l	0.050						
Chrysene	< 0.050		μg/l	0.050						
Dibenzo (a,h) anthracene	< 0.050		μg/l	0.050						
Fluoranthene	< 0.050		μg/l	0.050						
Fluorene	< 0.050		μg/l	0.050						
Indeno (1,2,3-cd) pyrene	< 0.050		μg/l	0.050						
Naphthalene	< 0.050		μg/l	0.050						
Phenanthrene	< 0.050		μg/l	0.050						
Pyrene	< 0.050		μg/l	0.050						
Surrogate: Benzo (e) pyrene-d12	1.10		μg/l		1.00		110	30-130		
LCS (1621185-BS2)					Pre	epared: 05-	Dec-16 An	alyzed: 06-D	ec-16	
Acenaphthene	0.706		μg/l	0.050	1.00		71	40-140		
Acenaphthylene	0.712		μg/l	0.050	1.00		71	40-140		
Anthracene	0.741		μg/l	0.050	1.00		74	40-140		
Benzo (a) anthracene	0.746		μg/l	0.050	1.00		75	40-140		
Benzo (a) pyrene	0.781		μg/l	0.050	1.00		78	40-140		
Benzo (b) fluoranthene	0.770		μg/l	0.050	1.00		77	40-140		
Benzo (g,h,i) perylene	0.756		μg/l	0.050	1.00		76	40-140		
Benzo (k) fluoranthene	0.793		μg/l	0.050	1.00		79	40-140		
Chrysene	0.756		μg/l	0.050	1.00		76	40-140		
Dibenzo (a,h) anthracene	0.774		μg/l	0.050	1.00		77	40-140		
Fluoranthene	0.733		μg/l	0.050	1.00		73	40-140		
Fluorene	0.732		μg/l	0.050	1.00		73	40-140		
Indeno (1,2,3-cd) pyrene	0.838		μg/l	0.050	1.00		84	40-140		
Naphthalene	0.758		μg/l	0.050	1.00		76	40-140		
Phenanthrene	0.714		μg/l	0.050	1.00		71	40-140		
Pyrene	0.750		μg/l	0.050	1.00		75	40-140		
Surrogate: Benzo (e) pyrene-d12	1.68		μg/l		2.00		84	30-130		
LCS Dup (1621185-BSD2)					Pre	epared: 05-	Dec-16 An	alyzed: 06-D	ec-16	
Acenaphthene	0.606		μg/l	0.050	1.00		61	40-140	15	20
Acenaphthylene	0.602		μg/l	0.050	1.00		60	40-140	17	20
Anthracene	0.595	QR2	μg/l	0.050	1.00		60	40-140	22	20
Benzo (a) anthracene	0.643		μg/l	0.050	1.00		64	40-140	15	20
Benzo (a) pyrene	0.686		μg/l	0.050	1.00		69	40-140	13	20
Benzo (b) fluoranthene	0.672		μg/l	0.050	1.00		67	40-140	14	20
Benzo (g,h,i) perylene	0.667		μg/l	0.050	1.00		67	40-140	13	20
Benzo (k) fluoranthene	0.700		μg/l	0.050	1.00		70	40-140	12	20
Chrysene	0.646		μg/l	0.050	1.00		65	40-140	16	20
Dibenzo (a,h) anthracene	0.653		μg/l	0.050	1.00		65	40-140	17	20
Fluoranthene	0.605		μg/l	0.050	1.00		60	40-140	19	20
Fluorene	0.625		μg/l	0.050	1.00		62	40-140	16	20
Indeno (1,2,3-cd) pyrene	0.740		μg/l	0.050	1.00		74	40-140	12	20
Naphthalene	0.630		μg/l	0.050	1.00		63	40-140	18	20
Phenanthrene	0.560	QR2	μg/l	0.050	1.00		56	40-140	24	20

# Semivolatile Organic Compounds by GCMS - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Batch 1621185 - SW846 3510C										
LCS Dup (1621185-BSD2)					Pre	epared: 05-E	Dec-16 Ar	nalyzed: 06-D	ec-16	
Pyrene	0.614		μg/l	0.050	1.00		61	40-140	20	20
Surrogate: Benzo (e) pyrene-d12	1.06		µq/l		1.00		106	30-130		

# **General Chemistry Parameters - Quality Control**

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Batch 1621045 - General Preparation										
Duplicate (1621045-DUP1)			Source: SC	<u> 29001-01</u>	Pre	epared & Ar	nalyzed: 01	-Dec-16		
pH	6.75		pH Units			6.74			0.1	5
Reference (1621045-SRM1)					<u>Pre</u>	epared & Ar	nalyzed: 01	-Dec-16		
рН	6.03		pH Units		6.00		100	97.5-102. 5		
Reference (1621045-SRM2)					<u>Pre</u>	epared & Ar	nalyzed: 01	-Dec-16		
рН	5.98		pH Units		6.00		100	97.5-102. 5		
Batch 1621065 - General Preparation										
Blank (1621065-BLK1)					Pre	epared: 02-	Dec-16 Ar	nalyzed: 06-D	ec-16	
Total Suspended Solids	< 0.5		mg/l	0.5						
LCS (1621065-BS1)					Pre	epared: 02-	Dec-16 Ar	nalyzed: 06-D	ec-16	
Total Suspended Solids	100		mg/l	10.0	100		100	90-110		

# **Subcontracted Analyses - Quality Control**

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Batch 369621A - 369621										
BLK (BV93731-BLK)					Pre	epared & Ar	nalyzed: 12-	-Dec-16		
Oil and Grease by EPA 1664A	< 1.4		mg/L	1.4	40			-		
LCS (BV93731-LCS)					Pre	epared: A	nalyzed: 12	-Dec-16		
Oil and Grease by EPA 1664A	39.70		ma/L	1.4	40		99	85-115		20

### **Notes and Definitions**

QR2 The RPD result exceeded the QC control limits; however, both percent recoveries were acceptable. Sample results for the QC batch were accepted based on percent recoveries and completeness of QC data.

dry Sample results reported on a dry weight basis

NR Not Reported

RPD Relative Percent Difference

OG The required Matrix Spike and Matrix Spike Duplicate (MS/MSD) for Oil & Grease method 1664B can only be analyzed when the client has submitted sufficient sample volume. An extra liter per MS/MSD is required to fulfill the method QC criteria. Please refer to Chain of Custody and QC Summary (MS/MSD) of the Laboratory Report to verify ample sample volume was submitted to fulfill the requirement.

pH The method for pH does not stipulate a specific holding time other than to state that the samples should be analyzed as soon as possible. For aqueous samples the 40 CFR 136 specifies a holding time of 15 minutes from sampling to analysis. Therefore all aqueous pH samples not analyzed in the field are considered out of hold time at the time of sample receipt. All soil samples are analyzed as soon as possible after sample receipt.

<u>Laboratory Control Sample (LCS)</u>: A known matrix spiked with compound(s) representative of the target analytes, which is used to document laboratory performance.

Matrix Duplicate: An intra-laboratory split sample which is used to document the precision of a method in a given sample matrix.

<u>Matrix Spike</u>: An aliquot of a sample spiked with a known concentration of target analyte(s). The spiking occurs prior to sample preparation and analysis. A matrix spike is used to document the bias of a method in a given sample matrix.

<u>Method Blank</u>: An analyte-free matrix to which all reagents are added in the same volumes or proportions as used in sample processing. The method blank should be carried through the complete sample preparation and analytical procedure. The method blank is used to document contamination resulting from the analytical process.

Method Detection Limit (MDL): The minimum concentration of a substance that can be measured and reported with 99% confidence that the analyte concentration is greater than zero and is determined from analysis of a sample in a given matrix type containing the analyte.

Reportable Detection Limit (RDL): The lowest concentration that can be reliably achieved within specified limits of precision and accuracy during routine laboratory operating conditions. For many analytes the RDL analyte concentration is selected as the lowest non-zero standard in the calibration curve. While the RDL is approximately 5 to 10 times the MDL, the RDL for each sample takes into account the sample volume/weight, extract/digestate volume, cleanup procedures and, if applicable, dry weight correction. Sample RDLs are highly matrix-dependent.

<u>Surrogate</u>: An organic compound which is similar to the target analyte(s) in chemical composition and behavior in the analytical process, but which is not normally found in environmental samples. These compounds are spiked into all blanks, standards, and samples prior to analysis. Percent recoveries are calculated for each surrogate.

<u>Continuing Calibration Verification:</u> The calibration relationship established during the initial calibration must be verified at periodic intervals. Concentrations, intervals, and criteria are method specific.

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☑ Standard TAT - 7 to 10 business days Special Handling:

Fediuring OCV	PECTRUM ANALYTICAL, INC.		

# CHAIN OF CUSTODY RECORD

	CHAIN OF CUSIODY RECOR	☐ Rush TAT - Date Needed:
SPECTRUM ANALYTICAL, INC. Featuring HANIBAL TECHNOLOGY	Page1 of1_	All TATs subject to laboratory approval Min. 24-hr notification needed for rushes Samples disposed after 60 days unless otherwise instructed.
Report To: Andrew Adams	Invoice To: Christopher Gill	Project No. Gulf Chelsea
Gulf Oil LP	Gulf Oil LP	
281 Eastern Ave	80 William St, Suite 400	Site Name: Guil Crieisea i eminial
Chelsea, MA 02150	Wellesley, MA 02481-3705	Location: 281 Eastern Ave, Chelsea State: MA
Telephone #: 617.884.5980	i.	Sampler(s): Andrew Adams
Project Mgr. Andrew Adams	P.O No.: Quote/RQN:	
- 1	4=HNO <sub>3</sub> 5=NaOH 6=Ascorbic Acid	List Preservative Code below: OA/OC Reporting Notes:
7=CH3OH 8=NaHSO <sub>4</sub> 9=Deionized Water 10=H <sub>3</sub> PO <sub>4</sub>	11= none 12= 41 2	2 11
	11 3	2 11

Condition upon receipt Custody Seals:  CAmbient Cleed Refrigerated	Greenbar Frankr	2 1/1/0 S E-mail to: aadams@gulfoil.com, cgill@gulfoil.com	Time: Temp °C					× ×	×	×	1 ×	naphthale	nzene & ne)	Containers : Analysis	11 3 2 11	List Preservative Code below:	
	101	Minus 12/11	Received by: Date:			- A	*	3 11.30-16 0760 6 SW	3 11-30-12 6900 6 SW 3	3 11.30 16 0700 G SW	3 11-36-16 6700 G SW	D: C=Compsite Time: Type Matrix # of VOA	A=Indoor/Ambient Air SG=Soil Gas  X2=X3=	sw=Surface Water ww=Waste Water	No.	1 3=H <sub>2</sub> SO <sub>4</sub> 4=HNO <sub>3</sub> 5=NaOH 6=Ascorbic Acid ater 10=H <sub>3</sub> PO <sub>4</sub> 11= none 12=	
N	Muny	awike	Relinquished by:					Outfall 003	Outfall 003	Outfall 003	29001 outfall 003	G= Grab  Lab ID: Sample ID:	O=Oil SO=Soil SL=Sludge A=I  X1=	<b>DW</b> =Dinking Water <b>GW</b> =Groundwater		F=Field Filtered 1=Na <sub>2</sub> S2O <sub>3</sub> 2=HCl 3=H <sub>2</sub> SO <sub>4</sub> 7=CH3OH 8=NaHSO <sub>4</sub> 9=Deionized Water 10=H <sub>3</sub> PO <sub>4</sub>	